

# 1-Decanamine, mono-DMTBS

**Inchi:** InChI=1S/C16H37NSi/c1-7-8-9-10-11-12-13-14-15-17-18(5,6)16(2,3)4/h17H,7-15H2,1-6H3  
**InchiKey:** QIJCOFUVDQSOI-UHFFFAOYSA-N  
**Formula:** C16H37NSi  
**SMILES:** CCCCCCCCCN[Si](C)(C)C(C)(C)C  
**Mol. weight [g/mol]:** 271.56

## Physical Properties

Property code	Value	Unit	Source
log10ws	-3.76		Crippen Method
logp	5.722		Crippen Method
rinpol	1647.00		NIST Webbook
rinpol	1647.00		NIST Webbook

## Sources

**NIST Webbook:** <http://webbook.nist.gov/cgi/cbook.cgi?ID=R64792&Units=SI>  
**Crippen Method:** <http://pubs.acs.org/doi/abs/10.1021/ci9903071>  
**Crippen Method:** [https://www.cheméo.com/doc/models/crippen\\_log10ws](https://www.cheméo.com/doc/models/crippen_log10ws)

## Legend

**log10ws:** Log10 of Water solubility in mol/l  
**logp:** Octanol/Water partition coefficient  
**rinpol:** Non-polar retention indices

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